

LUBRIZOL
INTER-OFFICE MEMORANDUM

(1)

To: GWBA

From: LAWM Date: _____

Subject: Level II Support - MoDTC

Cc: JACK, RVG, EAB, BRD, WBC, WMN, BDD

Gordon,

Attached is the Level II Support for 500 ppm Mo on top of [REDACTED]. Please forward this data to Steve and Mike at GM.

[REDACTED] is Lubrizol's general sales [REDACTED] product. The VM is Lubrizol's standard OCP. The base oils are Group 1 type.

Please note: We ran the IIIE on 250 ppm Mo since at 500 ppm Mo we failed on ORLD twice, 3.09 and 3.39. All formulas were exactly the same except for the level of MoDTC. The MoDTC used was Sakura-Lube.


Lew

LAWM/mjt
Attachment

**LEVEL 2 SUPPORT DATA FOR THE ADDITION OF MOLYBDENUM
DITHIOCARBAMATE**

Formulation	Baseline	Moly @ 500 ppm
Viscosity Grade	5W-30	5W-30
LZ 9800TM	X	X
Base Stocks		
Ashland 100N	88.0	88.0
Ashland 325N	12.0	12.0
Viscosity Modifier		
LZ 7070	8.20	8.20
Pour Point Depressant		
LZ 7749B	0.20	0.20
Dispersant/Inhibitor Package		
TBN	5.50	5.51
%Mo		0.044
%Mg	0.023	0.023
%Na	0.024	0.024
%Ca	0.092	0.092
%SA	0.659	0.659
%P	0.092	0.092
%Zn	0.101	0.101
%N	0.064	0.065
%S	0.485	0.460
CRC L-38	Pass	Pass
BWL (40 mg max.)	17.2	15.4
PSV (9.0 min.)	9.80	9.80
Sequence IID	Pass	Pass
AER (8.46 min.)	8.52	8.65
Sequence IIIE	Pass	Pass*
Hours to 375% Vis. Inc. (64 min.)	73.2	73.5
AES (9.16 min.)	9.58	9.56
APV (8.86 min.)	9.03	9.26
RLD (3.46 min.)	5.47	3.53
CLWM (64 um max.)	49	12
CLWA (30 um max.)	12.9	7.0
Oil Related Stuck Rings (none)	None	None
Sequence VE	Pass	Pass
AES (8.96 min.)	9.36	9.29
RCS (6.96 min.)	8.64	9.16
AEV (4.96 min.)	5.86	5.43
APV (6.46 min.)	6.82	7.39
CWM (15 um max.)	0.80	8.5
CWA (5 um max.)	1.03	3.87

* = % molybdenum = ~250 ppm.

**LEVEL 4 SUPPORT DATA FOR THE ADDITION OF MOLYBDENUM
DITHIOTOCARBAMATE**

Formulation	Baseline	Moly @ 500 ppm	Baseline	Moly @ 500 ppm
Viscosity Grad	5W-30	5W-30	5W-30	5W-30
OS Number	115072	128465	1131-108549	685-40120
Base Stocks				
Ashland 100N	88.0	88.0	90.0	90.0
Ashland 325N	12.0	12.0	10.0	10.0
Viscosity Modifier				
LZ 7070	8.20	8.20	8.10	8.10
Pour Point Depressant				
Viscoplex 1-330	-	-	0.20	0.20
LZ 7749B	0.20	0.20	-	-
Dispersant/Inhibitor Package				
TBN	5.50	5.51	5.12	5.10
%Mo	-	0.044	-	0.046
%Mg	0.023	0.023	0.037	0.037
%Na	0.024	0.024	0.059	0.059
%Ca	0.092	0.092	0.048	0.048
%SA	0.659	0.659	0.688	0.688
%P	0.092	0.092	0.095	0.095
%Zn	0.101	0.101	0.105	0.105
%N	0.064	0.065	0.053	0.053
%S	0.485	0.460	0.485	0.505
CRC L-38	Pass	Pass		
BWL (40 mg max.)	17.2	15.4		
PSV (9.0 min.)	9.80	9.80		
Sequence IID	Pass	Pass		
AER (8.46 min.)	8.52	8.65		
Sequence IIIE	Pass	Pass*		
Hours to 375% Vis. Inc. (64 min.)	73.2	73.5		
AES (9.16 min.)	9.58	9.56		
APV (8.86 min.)	9.03	9.26		
RLD (3.46 min.)	5.47	3.53		
CLWM (64 um max.)	49	12		
CLWA (30 um max.)	12.9	7.0		
Oil Related Stuck Rings (none)	None	None		
Sequence VE	Pass	Pass		
AES (8.96 min.)	9.36	9.29		
RCS (6.96 min.)	8.64	9.16		
AEV (4.96 min.)	5.86	5.43		
APV (6.46 min.)	6.82	7.39		
CWM (15 um max.)	0.80	8.5		
CWA (5 um max.)	1.03	3.87		
Sequence VIA	Pass		Fail	Pass
EFEI (1.06 min. for 5W-30)	1.13		0.99	1.33

* = % molybdenum = ~250 ppm.

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(2)

WOQ1 THE LUBRIZOL CORPORATION
INQUIRY SCREEN
BLEND REQUEST/MATERIAL DEFINITION

Material Name S010-4514-94-15, OS Number . . . OS128465

Enterer . . . RKW	Entry Date	[REDACTED]	Record Type	S
Réquestor . . . MW	Calc Unit. . .	W	CM Type . . .	BL
Priority . . . A	Blend Amount	82400 G		
Need By Date	Amt to Stock	33281 G	Picklist Stat	PICK

Project . . . T944514	Market Class	A	Blender	SAAB
Subset . . . 15	QA Approved		Blend Date	[REDACTED]
CMA# . . . Y	Maj/Min Mods	00 A	RM-4514940015-A-1	

Invent Date	[REDACTED]	Inventor . . .	MW	Invent Dept	5240	WKF
Perf Grade						
Visgrade . . . 5W-30		Visgr Source	SAE			
Waste Track	R NSW	7		Release Flag	N	

LZ Product

Press ENTER to Continue

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P/N

RESEARCH TEST RESULT (BOARD)

DATE: 07/15/03

PRODUCT: S010-4514-94-15, TEST NO:
AS OF 1/1/2000, USE MCR. ONLY HISTORICAL DATA WILL BE AVAILABLE IN WOBC

(3)

REMARKS:

1. REPLACED DISTRIBUTION DURING BREAK-IN.

CMA NO: RM-4514940015-A-1-VE-1-LZ-247
CHEM MATL NAME: S010-4514-94-15, OS NO: OS128465
TEST NAME: W080.045.07 COMP DATE: [REDACTED] DESC: SEQ.VE DEVL.
RUN NUMBER: 1 LOCATION: WICKLIFFE
VALIDITY: VALID QUALIFIED: QUALIFIED

ID NUMBER: 247-5-41 FUEL BATCH: 43
TEST HOUR: 288.0 ENGINE NO: 81-94-
PASS/FAIL: PASS

SLUDGE DEPOSITS	RATING	RESULT	ADJUSTED	VARNISH DEPOSITS	RATING	RESULT	ADJUSTED
ROCKER ARM COVER	9.16	9.16	-----	PISTON SKIRTS	7.39	7.39	-----

ENTER - NEXT PAGE, P/-1 - PREVIOUS PAGE, CLEAR - EXIT TRANSACTION

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0Q5

THE LUBRIZOL CORPORATION
INQUIRY SCREEN
FORMULA AT ACTUAL STATE

07/18/03
15:42:38

Material Name	Calc Unit . .	CM Type	BL	OS Number
S010-4514-94-15,	W			OS128465
265-X6052,P	BASE STOCKS			
265-X6047,H				
S762-4513-94-134,S				
S762-4850-94-718,B				
S010-4514-94-14,- CONC				
		Pct/PPM	Use	Pct W Pct V OS#
		88.0000	B	71.1920 72.0649 OS75287Y
		12.0000	B	9.7080 9.6219 OS75863K
		8.2000	O	8.2000 8.2866 OS115849L
		0.2000	O	0.2000 0.1922 OS118095C
		10.7000	O	10.7000 9.8344

Press ENTER to continue

WOQ5MSGIN Actual information displayed successfully.

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NOQ6

THE LUBRIZOL CORPORATION
INQUIRY SCREEN
THEORETICAL ANALYSIS

07/18/03
15:41:55

Material Name 265-X6052,P OS Number OS75287Y
Waste Tracking Number R NSW 7
Ela Name . . . 265-X6052

%Oil	100.0000	%Ca	%S	.2000	%Mo
Sp.Gr	.8644	%Mn	%N	.0020	%Pb
TBN		%Co	%B		%K
TAN		%Mg	%Cu		%Na
%SA		%Ba	%Cl		%Sr
%Zn		%P	%Li		%Si

Title Title Name . . . 265-X6052
* (MT) VALVOLINE/ASHLAND CATLETTSBURG 100N HT *USE DIL OIL

Press ENTER to continue
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WOQ6

THE LUBRIZOL CORPORATION
INQUIRY SCREEN
THEORETICAL ANALYSIS

07/18/03
10:09:46

Material Name 265-X6047,H
Waste Tracking Number R NSW 7
Ela Name . . . 265-X6047

OS Number OS75863K

%Oil	100.0000	%Ca	%S	.3900	%Mo
Sp.Gr	.8820	%Mn	%N	.0040	%Pb
TBN		%Co	%B		%K
TAN		%Mg	%Cu		%Na
%SA		%Ba	%Cl		%Sr
%Zn		%P	%Li		%Si

Title Title Name . . 265-X6047,H
VALVOLINE, INC.

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WOQ5

THE LUBRIZOL CORPORATION
INQUIRY SCREEN
FORMULA AT ACTUAL STATE

07/18/03
15:43:18

Material Name S010-4514-94-14,

Calc Unit . . W CM Type CN

Ingredient	Pct/PPM	Use	Pct W	Pct V	OS#
S015-4041-94-17,J	9.5000	O	88.7850	87.8868	OS115020C
S217-8226-95-01,C	1.2000	O	11.2150	12.1132	

Press ENTER to continue

WOQ5MSGIN Actual information displayed successfully.

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WOQ3

THE LUBRIZOL CORPORATION
INQUIRY SCREEN
RECEIVE OUTSIDE MATERIAL

07/18/03
15:44:30

Material Name	S217-8226-95-01, ←	OS Number	OS117924	
Enterer . . .	KJL	Requestor . . .		
Project . . .	T958226	Record Type . . .	S	
Subset . . .		CM Type . . .	UN	
CMA # . . .		Invent Date . . .		
Receipt Date		Document Type		
Receipt Amount	1.000	U/M G	Document Number	
Supplier . . .	TONEN		Inv Amount	1
Supp Batch Id				
Visgrade . . .		Visgr Source	Waste Track R	NSW
SpGr8800	MSDS Avail ?	Competitive Mat	7
Blend Title	MO-DTC SAKURA LUBE #100 ←			

Comment . . . TONEN

LZ Material

Press ENTER to Continue

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WOQ6

THE LUBRIZOL CORPORATION
INQUIRY SCREEN
THEORETICAL ANALYSIS

07/18/03
15:44:46

Material Name S217-8226-95-01,
Waste Tracking Number R NSW 7
Ela Name . . . S217-8226-95-01

%Oil	80.0000	%Ca	%S	4.1600	%Mo	4.1500
Sp.Gr	.8800	%Mn	%N		%Pb	
TBN	1.7000	%Co	%B		%K	
TAN		%Mg	%Cu		%Na	.0090
%SA	6.2528	%Ba	%Cl		%Sr	
%Zn		%P	.0160	%Li		%Si

Title Title Name . . . S217-8226-95-01,
TONEN

Press ENTER to continue
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WOQ5

THE LUBRIZOL CORPORATION
INQUIRY SCREEN
FORMULA AT ACTUAL STATE07/18/03
15:45:19

Material Name	S015-4041-94-17,	OS Number	OS115020
Calc Unit . .	W CM Type CN	Pct	V OS#
Ingredient	Pct/PPM	Use	
-C	4.0300	O	42.4210 44.5216 OS48147K
	0.8600	O	9.0526 10.0707
	1.3700	O	14.4211 14.8903 OS28802Y
-d	0.9200	O	9.6842 7.7467
- e-1	0.6800	O	7.1579 7.3149 OS128676
- e-2	0.3000	O	3.1579 3.1944 OS47056H
	0.5100	O	5.3684 4.7312
	0.3800	O	4.0000 3.3994 OS45933AA
	0.2000	O	2.1053 1.8676 OS101888A
	0.2500	O	2.6316 2.2632 OS84147F
	90.0000	T	0.0947 0.1101

Press ENTER to continue

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C ↗ [REDACTED]

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Intermediate Chemical Information

Last Updated: 23JUL03

Generic Name: Polyolefin amide alkeneamine**MSDS Chemical Name(s):** Polyisobutenylsuccinic anhydride, product with polyethyleneamines**Primary End Use(s):** DISPERSANT**Secondary End Use(s):** VISCOSITY INDEX IMPROVER**LZ Product Numbers:** [REDACTED]**Waste Tracking Number:** R-NSW-7**Functional Structures**

- SUCCINIMIDE
- SUCCINIMIDE, FROM CHLORINE SUCCAN
- SUCCINIMIDE, CO:N(1/0.83)=(6/5) EG [REDACTED]

Backbone Structures

- ALIPHATIC
- BRANCHED
- MOLECULAR WT SIDE CHAIN 1601-3000 (PBU)
- MOLECULAR WT SIDE CHAIN NUMBER AVERAGE
- OLEFINIC (DIENE, ALKYNE, ETC)
- POLYMÉR
- POLYMER, BUTYLENES
- POLYMER, HOMOPOLYMER

Elementals - Theory

%B	0	%CO		%MG		%NA		%SI	
%BA		%CU		%MN		%P		%SR	
%CA		%K		%MO		%PB		%ZN	0
%CL	0.3	%LI		%N	0.9	%S	0		

Other Characteristics - Theory

%Sulfonate		Conv_Sulfonate		%Sulfated Ash	
%Salicylate				ZDP Primary %P	
%Phenate		Conv_Phenate		ZDP Secondary %P	
%Carboxylate		Conv_Carboxylate		ZDP Aromatic %P	
%Thiosulfate		Conv_Thiosulfate			
%Oil	55	Specific Gravity	0.91	TBN	15

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Intermediate Chemical Information

Last Updated: 23JUL03

Generic Name: Zinc alkylidithiophosphate**MSDS Chemical Name(s):** Zinc O,O-di(1-methylethyl 1,3 dimethylbutyl) dithiophosphate**Primary End Use(s):** ANTIOXIDANT, ANTIWEAR AGENT**LZ Product Numbers:** [REDACTED]**Waste Tracking Number:** R-CRCD-L**Functional Structures**

- CONV = 100-150 / NEUTRALS (EG [REDACTED])
- DITHIOPHOSPHATE (DTP)
- DITHIOPHOSPHATE, SALT *
- DITHIOPHOSPHATE, SECONDARY
- SALT, ZINC

Backbone Structures

- ALIPHATIC
- BRANCHED
- MOLECULAR WT SIDE CHAIN UNDER 169 (

Elementals - Theory

%B		%CO		%MG		%NA		%SI	
%BA		%CU		%MN		%P	10	%SR	
%CA		%K		%MO		%PB		%ZN	11.05
%CL	0.0001	%LI		%N		%S	21		

Other Characteristics - Theory

%Sulfonate		Conv_Sulfonate		%Sulfated Ash	16.575
%Salicylate				ZDP Primary %P	0
%Phenate		Conv_Phenate		ZDP Secondary %P	10
%Carboxylate		Conv_Carboxylate		ZDP Aromatic %P	0
%Thiosulfate		Conv_Thiosulfate			
%Oil	9	Specific Gravity	1.2	TBN	5

e-1 → [REDACTED]

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*Lubrizol Confidential***Obsolete Intermediate Chemical Information**

Last Updated: 23JUL03

Generic Name: Alkaryl amine**MSDS Chemical Name(s):** Dinonyldiphenylamine**Primary End Use(s):** ANTIOXIDANT**Performance Synonyms:** 1050-4367, S348-6531-86-12, A755-630-L-80, PRIOLUBE 3999, 1131-88924, OIU-38**Waste Tracking Number:** R-HZL DFLT-R**Functional Structures**

- AMINE, AROMATIC
- AMINE, MONO
- AMINE, SECONDARY

Backbone Structures

- AROMATIC
- BRANCHED
- INHIBITOR - AMINE
- MOLECULAR WT OVERALL 282-519 (>C20-C37)
- MOLECULAR WT SIDE CHAIN UNDER 169 (

Elementals - Theory

%B		%CO	%MG		%NA		%SI	
%BA		%CU	%MN		%P		%SR	
%CA		%K	%MO		%PB		%ZN	
%CL	0.015	%LI	%N	3.2	%S			

Other Characteristics - Theory

%Sulfonate		Conv_Sulfonate		%Sulfated Ash	
%Salicylate				ZDP Primary %P	
%Phenate		Conv_Phenate		ZDP Secondary %P	
%Carboxylate		Conv_Carboxylate		ZDP Aromatic %P	
%Thiosulfate		Conv_Thiosulfate			
%Oil	12	Specific Gravity	0.94	TBN	120

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Obsolete Intermediate Chemical
Information

Last Updated: 23JUL03

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Generic Name: Olefin sulfide**MSDS Chemical Name(s):** Alkyl(C12-C18) sulfide**Primary End Use(s):** ANTIOXIDANT**Secondary End Use(s):** ANTIWEAR AGENT**Waste Tracking Number:** R-NHW-1**Functional Structures**

- SULFIDE, UNKNOWN (EG SULFURIZED OLEFINS)

Backbone Structures

- ALIPHATIC
- BRANCHED
- INHIBITOR - SULFURIZED OLEFIN
- MOLECULAR WT SIDE CHAIN 169-225(C12-C16)

Elementals - Theory

%B		%CO		%MG		%NA		%SI	
%BA		%CU		%MN		%P		%SR	
%CA		%K		%MO		%PB		%ZN	
%CL	0.5	%LI		%N		%S	19		

Other Characteristics - Theory

%Sulfonate	Conv_Sulfonate		%Sulfated Ash	
%Salicylate			ZDP Primary %P	
%Phenate	Conv_Phenate		ZDP Secondary %P	
%Carboxylate	Conv_Carboxylate		ZDP Aromatic %P	
%Thiosulfate	Conv_Thiosulfate			
%Oil	15	Specific Gravity	0.95	TBN
				0